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IN THIS ISSUE

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Cover
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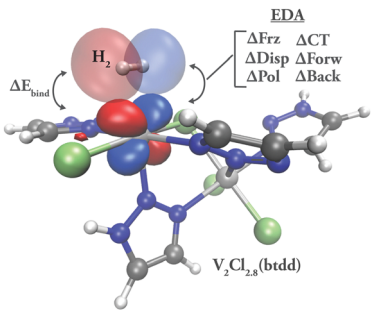

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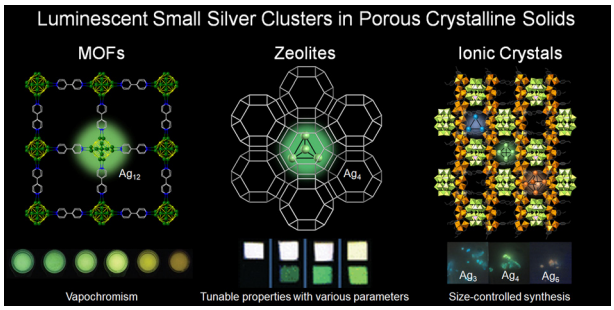

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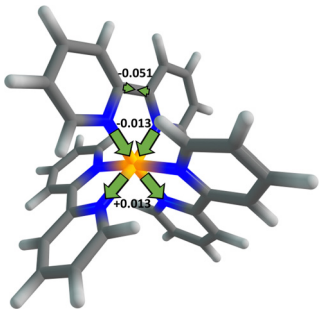

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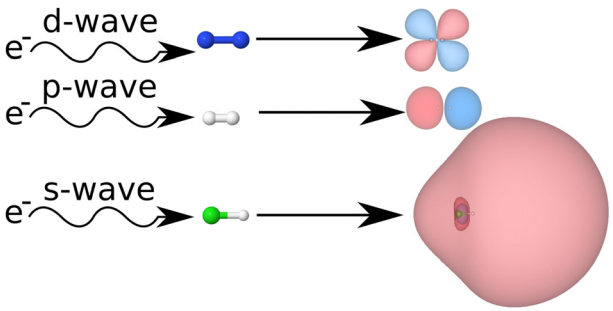

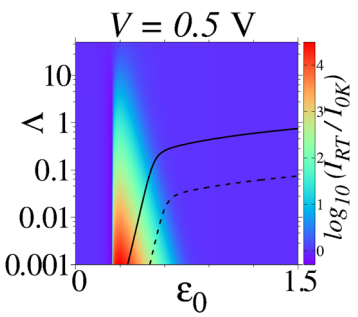

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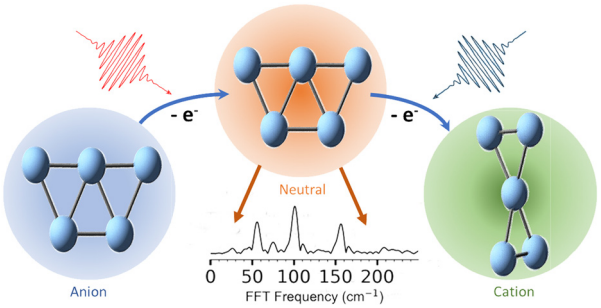

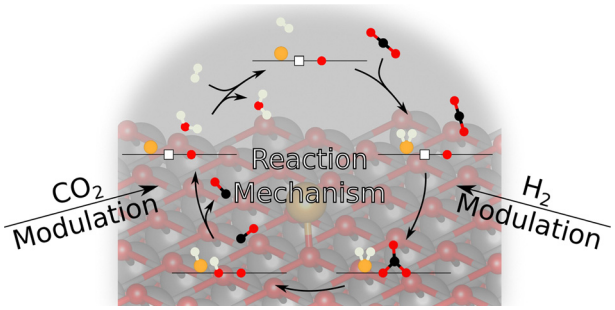

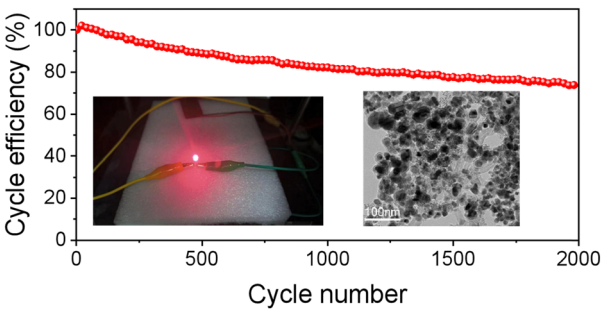

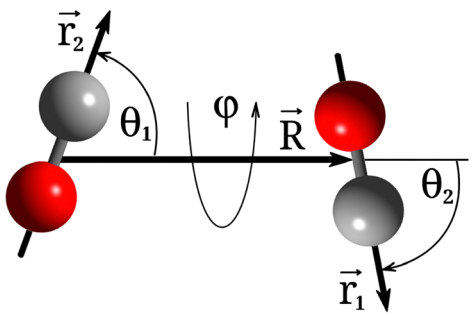

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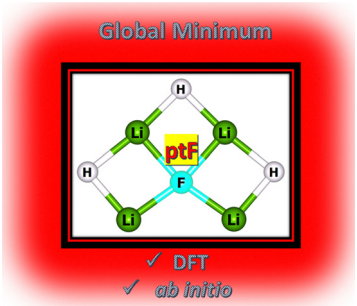

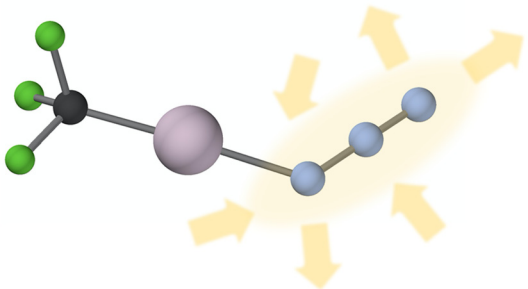

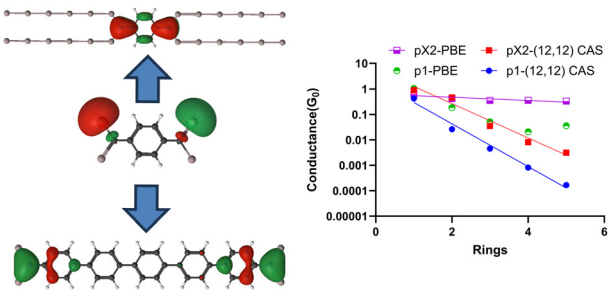

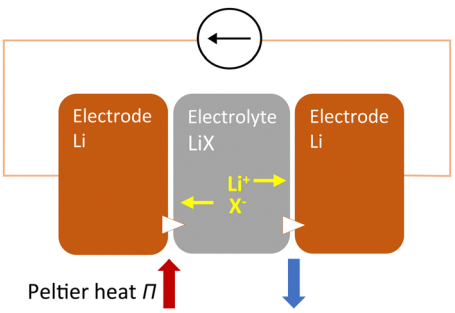

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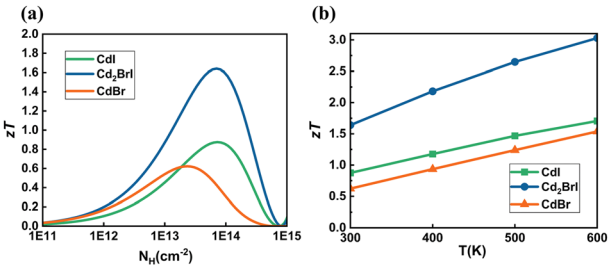

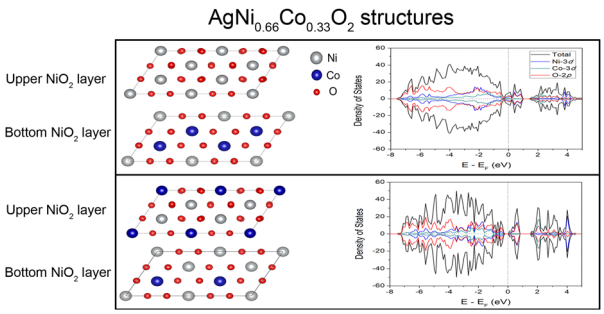

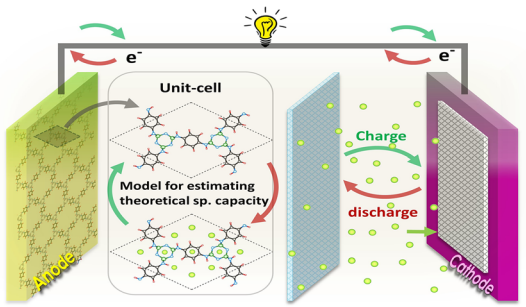

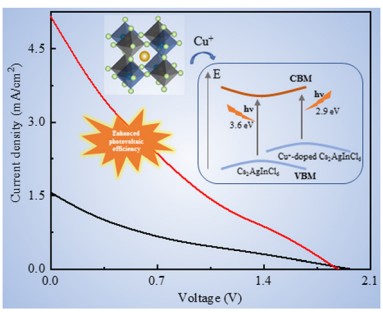

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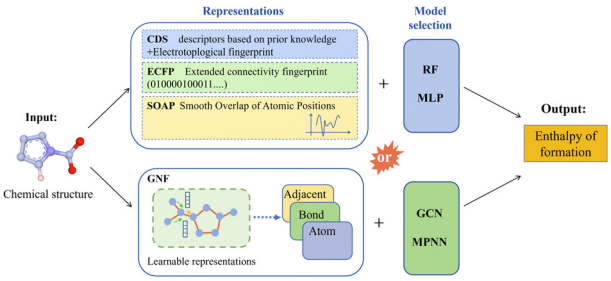

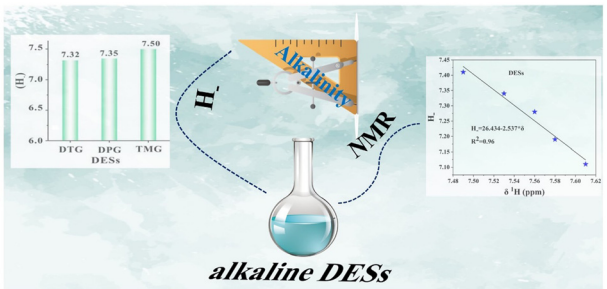

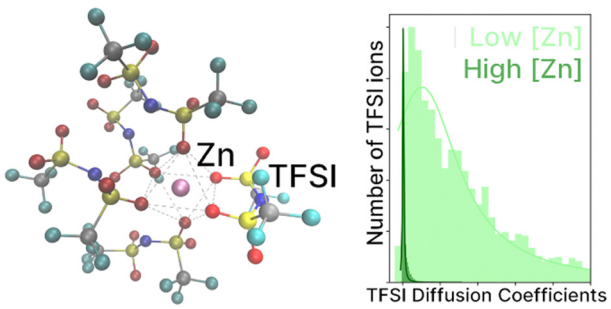

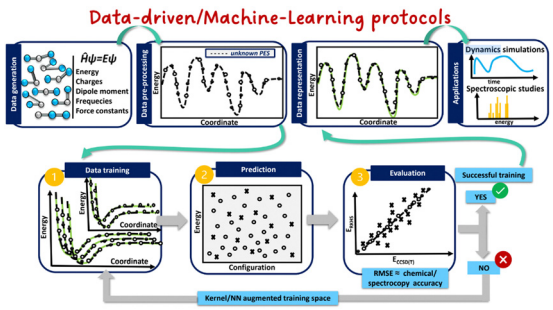

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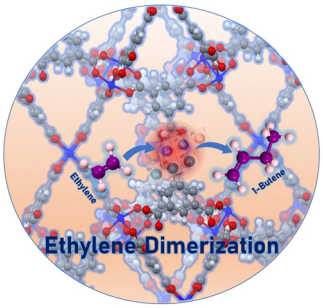

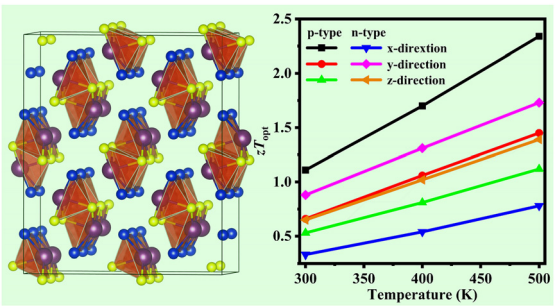

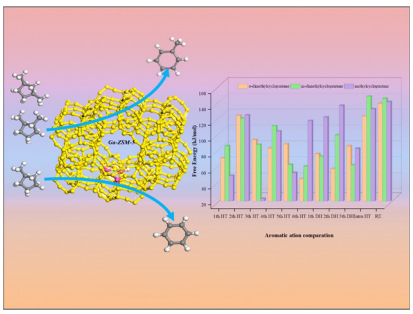

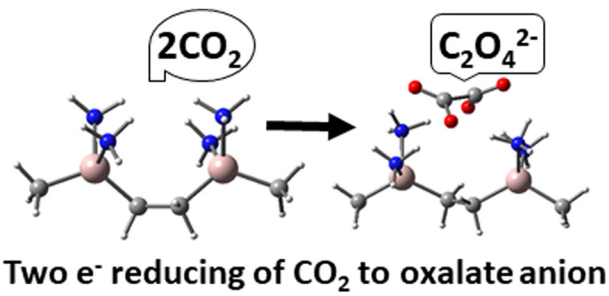

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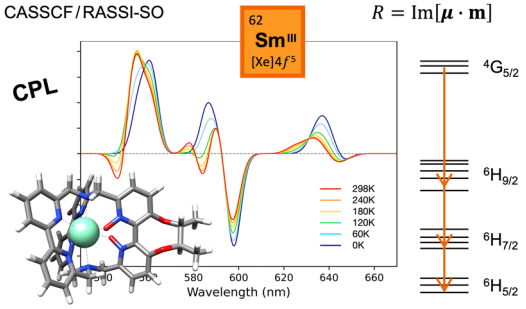
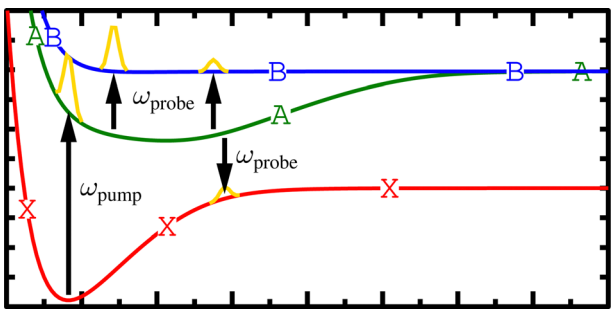
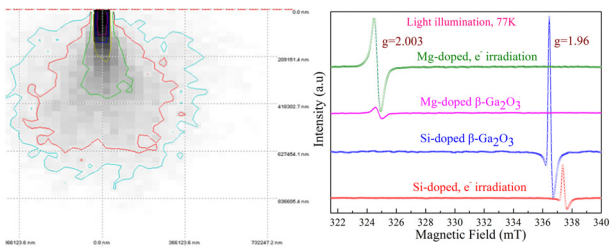
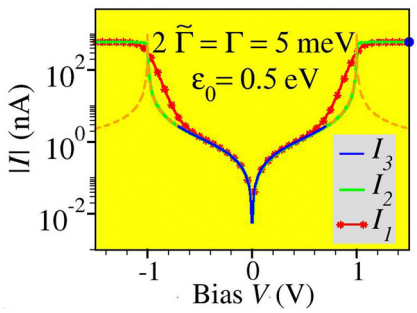






COMMENTS



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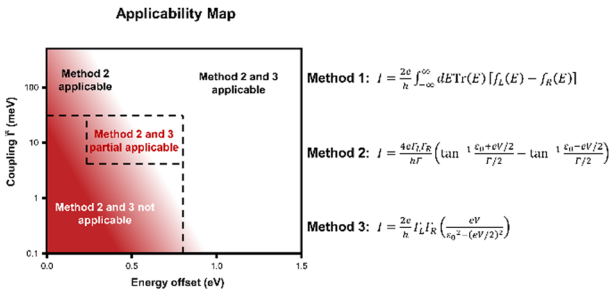

Reply to the 'Comment on "A single level tunneling model for molecular junctions: evaluating the simulation methods"' by I Baldea, *Phys. Chem. Chem. Phys.*, 2024, 26, D2CP05110A (http://D2CP05110A)

Zheyang Li and Xi Yu*

CORRECTION

Correction: Impact of temperature-dependent non-PAN peroxynitrate formation, RO_2NO_2, on nighttime atmospheric chemistry

Michelle Färber, Luc Vereecken, Hendrik Fuchs, Georgios I. Gkatzelis, Franz Rohrer, Sergej Wedel, Andreas Wahner and Anna Novelli*